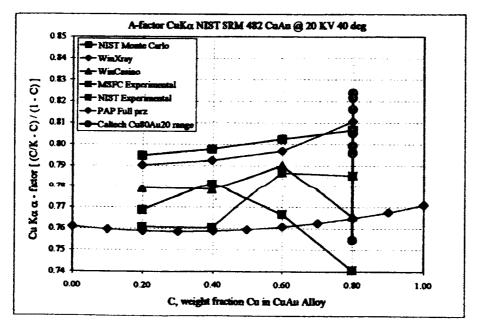
## A Comparison of Experimental EPMA Data and Monte Carlo Simulations

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Monte Carlo (MC) modeling shows excellent prospects for simulating electron scattering and x-ray emission from complex geometries, and can be compared to experimental measurements using electronprobe microanalysis (EPMA) and  $\phi(\rho z)$  correction algorithms. Experimental EPMA measurements made on NIST SRM 481 (AgAu) and 482 (CuAu) alloys, at a range of accelerating potential and instrument take-off angles, represent a formal microanalysis data set that has been used to develop  $\phi(\rho z)$ correction algorithms. The accuracy of MC calculations obtained using the NIST, WinCasino, WinXray, and Penelope MC packages will be evaluated relative to these experimental data. The αfactor method has previously been used to evaluate systematic errors in the analysis of semiconductor and silicate minerals [1,2,3], and will be used here to compare the accuracy of experimental and calculated x-ray data. X-ray intensities calculated by MC are used to generate \alpha-factors using the certificated compositions in the CuAu binary relative to pure Cu and Au standards. MC-generated x-ray intensities have a "built-in" atomic number correction, and are further corrected for absorption and characteristic and continuum fluorescence by φ(pz) correction algorithms. Preliminary results for αfactor analysis of Cu Ka in SRM 482 at 20 kV and 40 degree takeoff angle are shown in figure 1. For these data there is ~ 5% agreement between intensities calculated by MC and the PAP  $\phi(\rho z)$  algorithm compared to experimental EPMA data acquired from three different instruments. This excellent agreement indicates that MC modeling can be successfully used to calculate x-ray intensities for quantitative EPMA.



## References

- [1] P. Carpenter, Microbeam Analysis (2000), 433.
- [2] P. Carpenter, Microscopy and Microanalysis (1999), 74.
- [3] P. Carpenter, J. Armstrong, Microscopy and Microanalysis (1998), 224.